

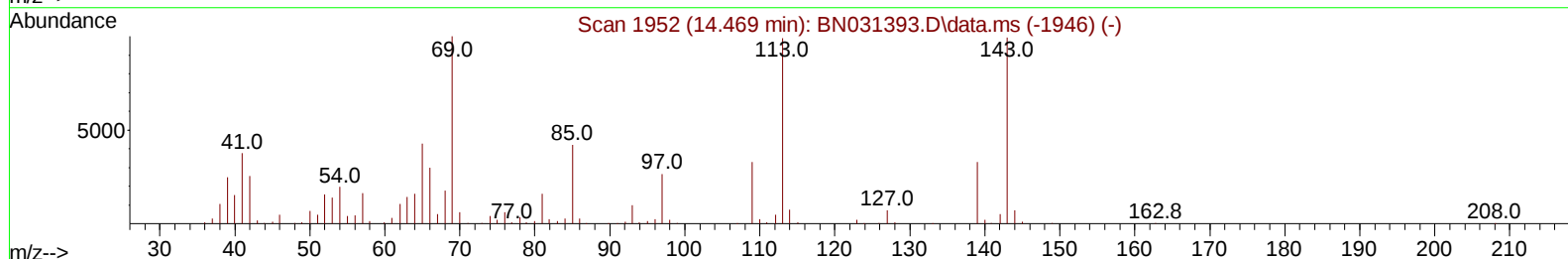
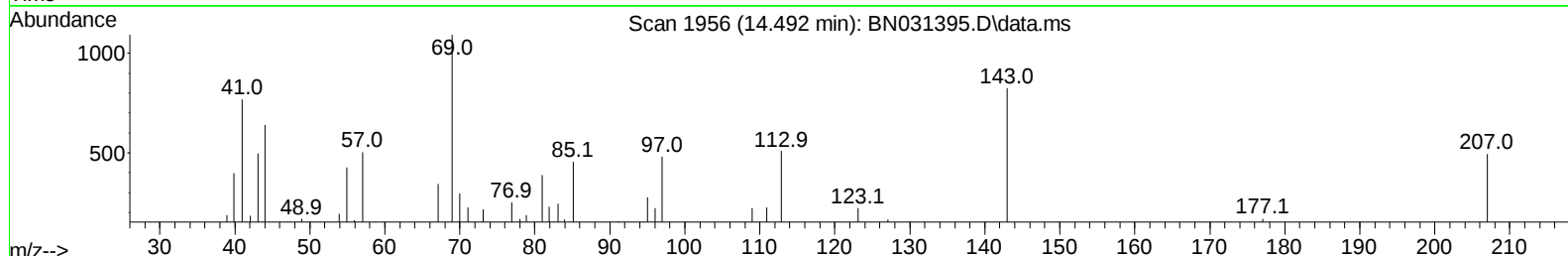
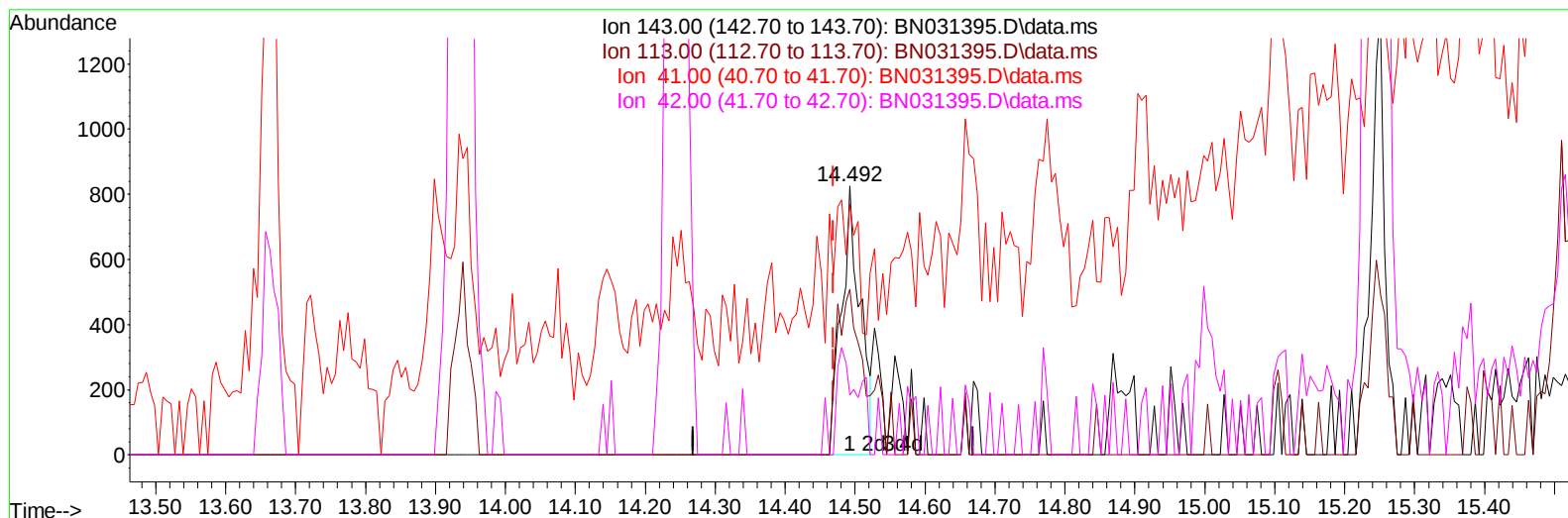
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN050624\
 Data File : BN031395.D
 Acq On : 06 May 2024 12:05
 Operator : MA/JU
 Sample : P2279-07
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 CORR0

Manual IntegrationsAPPROVED

Quant Time: May 06 12:42:36 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN050324.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri May 03 14:17:04 2024
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 05/07/2024
 Supervised By :mohammad ahmed 05/08/2024



TIC: BN031395.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.492min (+ 0.023) 0.45 ng/u1

response 1567

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.20	61.58#
41.00	39.50	93.21#
42.00	26.20	22.18

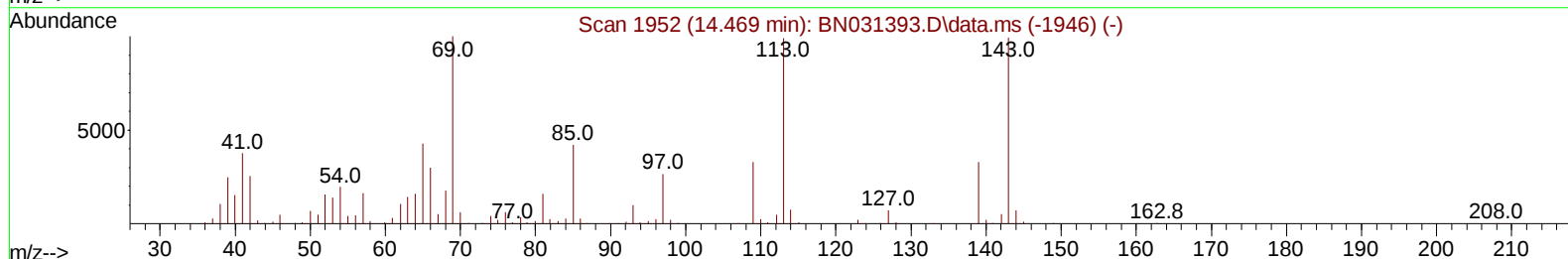
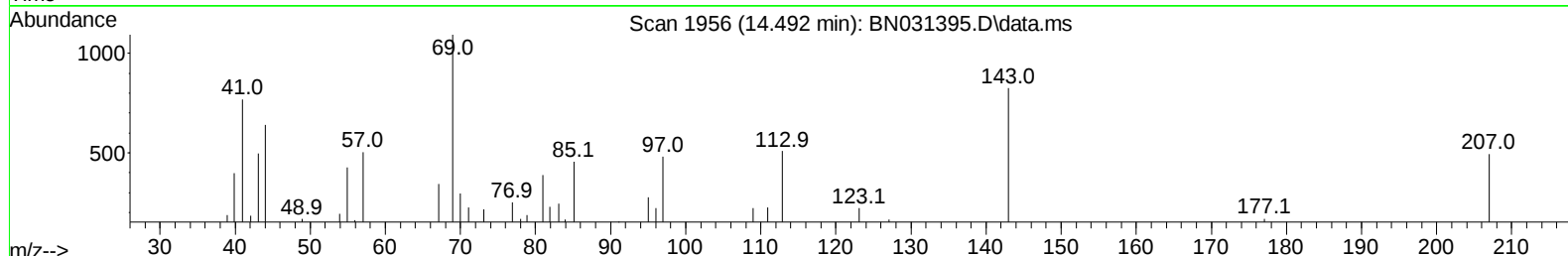
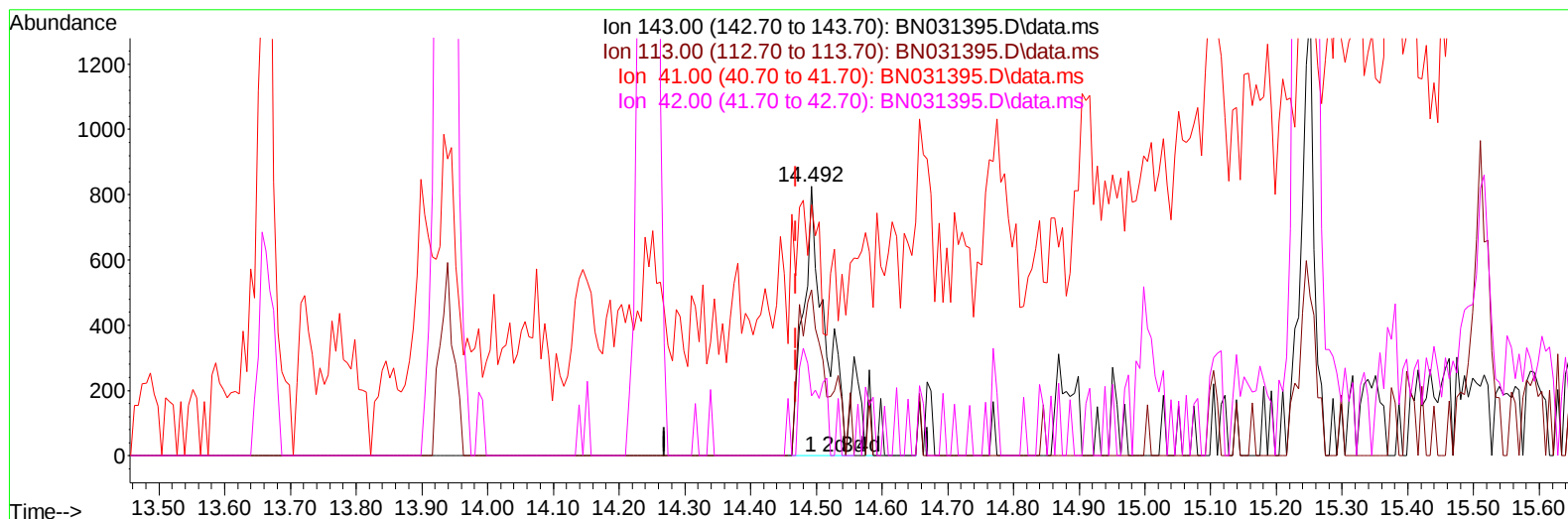
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Quant Time: May 06 12:42:36 2024
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TIC: BN031395.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.492min (+ 0.023) 0.66 ng/ul m

response 2335

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.20	61.58#
41.00	39.50	93.21#
42.00	26.20	22.18

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 Sample : P2279-07
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 CORR0

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 05/07/2024
 Supervised By :mohammad ahmed 05/08/2024

Quant Time: May 06 12:43:54 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN050324.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri May 03 14:17:04 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.593	152	99942	20.000	ng/ul	0.00
20) Naphthalene-d8	10.375	136	409704	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.245	164	275090	20.000	ng/ul	0.00
64) Phenanthrene-d10	16.998	188	578462	20.000	ng/ul	0.00
79) Chrysene-d12	21.233	240	542515	20.000	ng/ul	0.00
88) Perylene-d12	24.110	264	522136	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.117	96	13260	5.742	ng/uL	0.00
4) Pyridine-d5	3.522	84	79395	13.228	ng/ul	0.00
7) Phenol-d5	6.769	99	50348	6.607	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.940	67	149773	34.804	ng/ul	0.00
11) 2-Chlorophenol-d4	7.128	132	154131	24.253	ng/ul	0.00
15) 4-Methylphenol-d8	8.299	113	90831	14.374	ng/ul	-0.01
21) Nitrobenzene-d5	8.752	128	113480	36.763	ng/ul	0.00
24) 2-Nitrophenol-d4	9.463	143	49877	14.621	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.999	165	156535	25.531	ng/ul	0.00
31) 4-Chloroaniline-d4	10.516	131	198757	23.283	ng/ul	0.00
46) Dimethylphthalate-d6	13.663	166	788038	39.812	ng/ul	0.00
49) Acenaphthylene-d8	13.940	160	840999	38.606	ng/ul	0.00
54) 4-Nitrophenol-d4	14.492	143	2335m	0.664	ng/ul	0.02
60) Fluorene-d10	15.245	176	674705	41.707	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.375	200	7195	2.034	ng/ul	0.00
73) Anthracene-d10	17.098	188	1128646	44.803	ng/ul	0.00
81) Pyrene-d10	19.404	212	1367680	44.395	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.921	264	1199246	48.149	ng/ul	0.00
Target Compounds						
86) Bis(2-ethylhexyl)phtha...	21.151	149	49148	2.021	ng/ul	Qvalue 96

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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